



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 05:00 AM UTC

PDB ID : 4AOY / pdb_00004aoy
Title : Open CtIDH. The complex structures of Isocitrate dehydrogenase from Clostridium thermocellum and Desulfotalea psychrophila, support a new active site locking mechanism
Authors : Leiros, H.-K.S.; Fedoy, A.-E.; Leiros, I.; Steen, I.H.
Deposited on : 2012-03-30
Resolution : 2.35 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

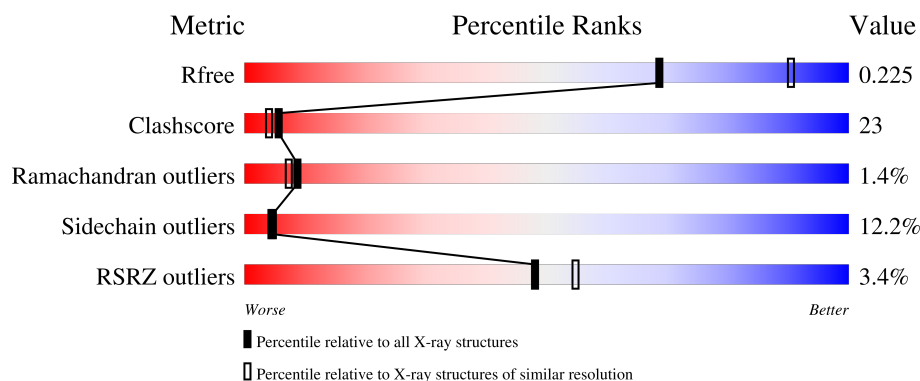
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	402	4% 59% 27% 6% 6%
1	B	402	4% 56% 27% 8% 7%
1	C	402	0% 66% 21% • 8%
1	D	402	3% 58% 28% 7% 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11678 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called ISOCITRATE DEHYDROGENASE [NADP].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	376	Total	C	N	O	S	0	0	0
			2917	1854	488	558	17			
1	B	372	Total	C	N	O	S	0	0	0
			2859	1817	474	552	16			
1	C	371	Total	C	N	O	S	0	0	0
			2884	1837	477	554	16			
1	D	375	Total	C	N	O	S	0	0	0
			2930	1869	485	560	16			

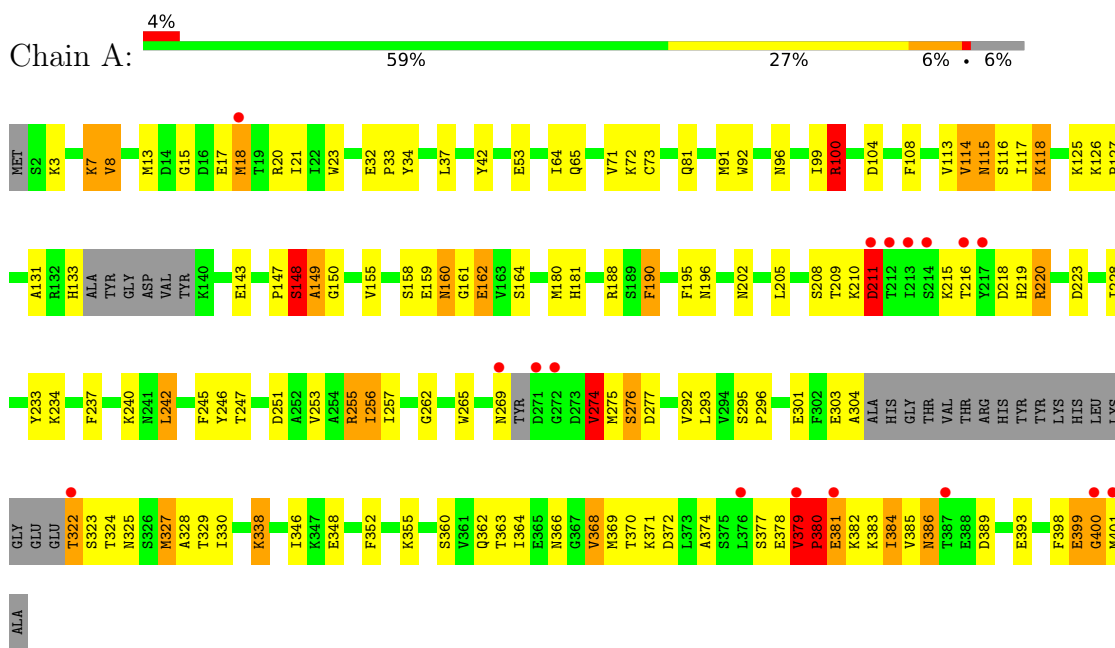
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	21	Total	O	0	0
			21	21		
2	B	15	Total	O	0	0
			15	15		
2	C	22	Total	O	0	0
			22	22		
2	D	30	Total	O	0	0
			30	30		

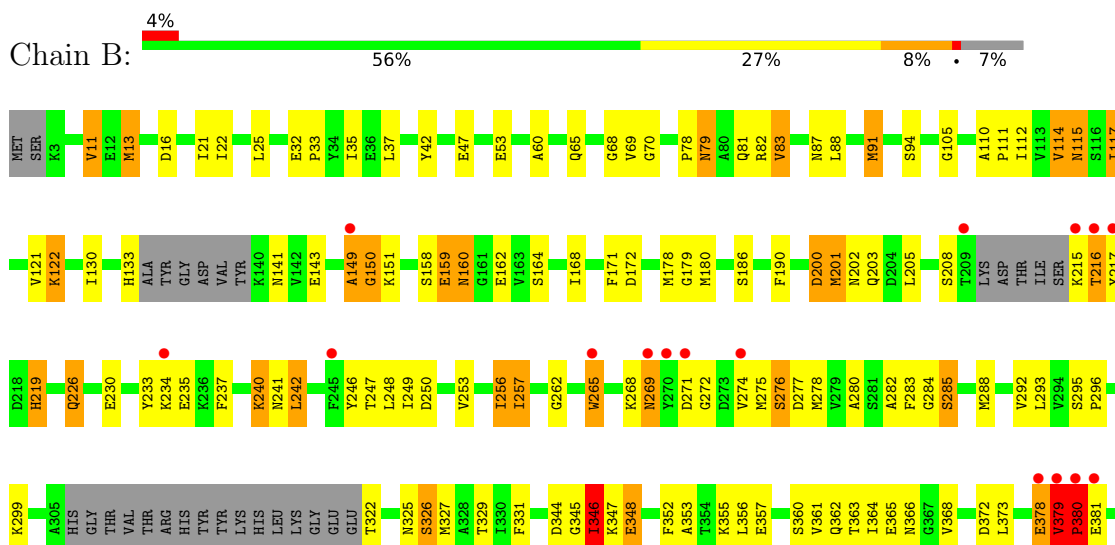
3 Residue-property plots

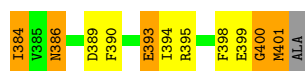
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: ISOCITRATE DEHYDROGENASE [NADP]

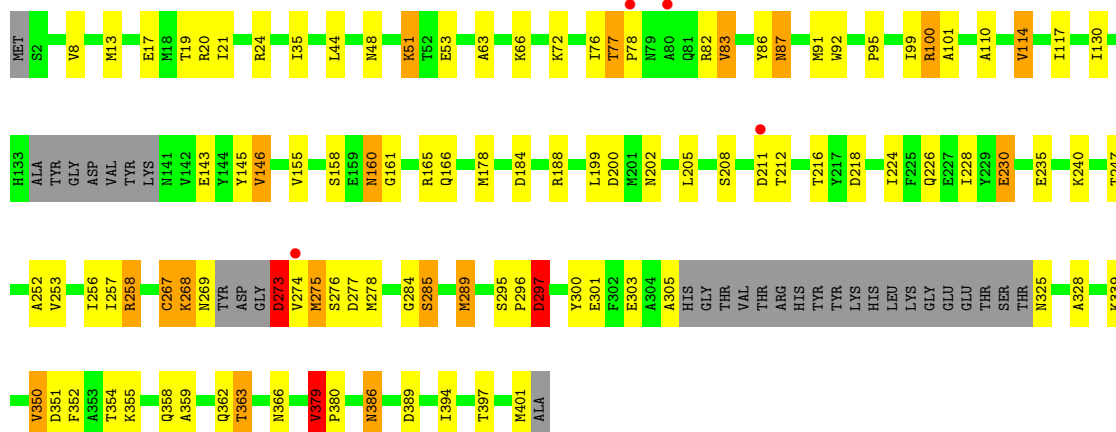


• Molecule 1: ISOCITRATE DEHYDROGENASE [NADP]

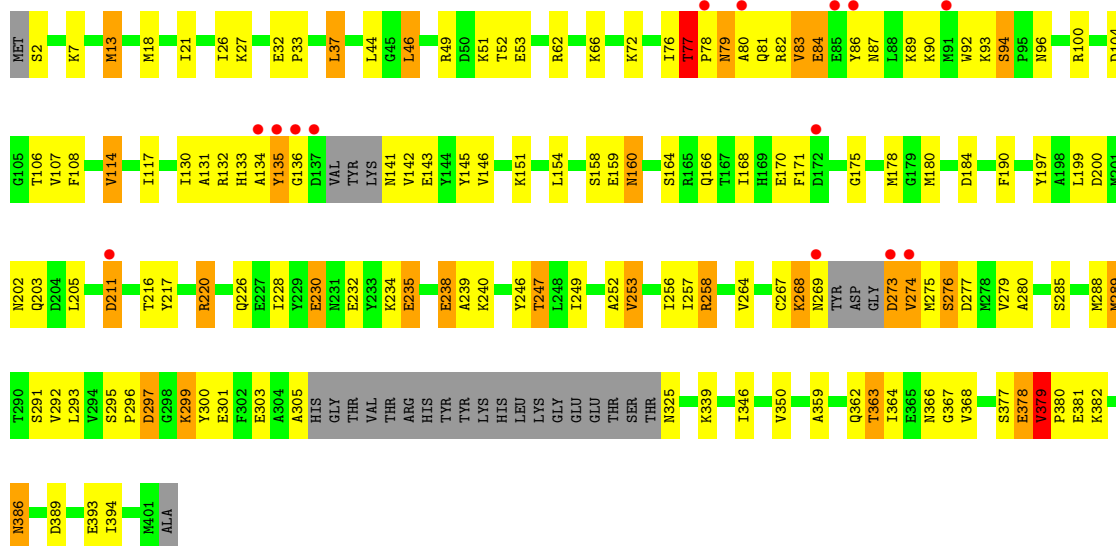




• Molecule 1: ISOCITRATE DEHYDROGENASE [NADP]



• Molecule 1: ISOCITRATE DEHYDROGENASE [NADP]



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.13Å 106.84Å 154.50Å 90.00° 93.63° 90.00°	Depositor
Resolution (Å)	20.00 – 2.35 20.00 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.7 (20.00-2.35) 99.6 (20.00-2.35)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.210 , 0.258 0.211 , 0.225	Depositor DCC
R_{free} test set	2280 reflections (3.01%)	wwPDB-VP
Wilson B-factor (Å ²)	40.4	Xtriage
Anisotropy	0.050	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11678	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.14% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.84	0/2970	1.15	12/4011 (0.3%)
1	B	0.83	0/2913	1.24	16/3936 (0.4%)
1	C	0.92	0/2938	1.14	18/3966 (0.5%)
1	D	0.92	0/2985	1.09	14/4029 (0.3%)
All	All	0.88	0/11806	1.15	60/15942 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	1
All	All	0	3

There are no bond length outliers.

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	379	VAL	CA-C-N	-22.05	92.28	119.84
1	B	379	VAL	C-N-CA	-22.05	92.28	119.84
1	A	379	VAL	CA-C-N	-13.08	103.50	119.84
1	A	379	VAL	C-N-CA	-13.08	103.50	119.84
1	B	149	ALA	N-CA-C	10.37	128.37	111.37
1	C	379	VAL	CA-C-N	10.30	130.61	119.28
1	C	379	VAL	C-N-CA	10.30	130.61	119.28
1	D	77	THR	CA-C-N	9.75	132.03	119.84
1	D	77	THR	C-N-CA	9.75	132.03	119.84
1	B	150	GLY	N-CA-C	8.13	132.45	113.18
1	C	379	VAL	N-CA-CB	7.20	117.41	109.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	379	VAL	CA-C-N	6.77	126.72	119.28
1	D	379	VAL	C-N-CA	6.77	126.72	119.28
1	A	380	PRO	CB-CA-C	-6.75	100.43	111.56
1	C	8	VAL	CA-C-N	-6.62	113.13	120.14
1	C	8	VAL	C-N-CA	-6.62	113.13	120.14
1	B	400	GLY	N-CA-C	-6.61	97.52	113.18
1	B	380	PRO	CB-CA-C	-6.46	100.91	111.56
1	A	274	VAL	CB-CA-C	-6.44	103.76	111.81
1	D	135	TYR	N-CA-C	6.44	120.09	107.98
1	A	150	GLY	N-CA-C	-6.33	101.39	110.38
1	B	151	LYS	N-CA-C	6.21	118.64	108.52
1	B	381	GLU	N-CA-C	-6.18	100.12	109.95
1	C	77	THR	CB-CA-C	5.93	116.61	109.85
1	A	190	PHE	N-CA-C	-5.92	104.75	111.14
1	A	211	ASP	N-CA-C	5.89	123.36	110.80
1	B	168	ILE	N-CA-C	-5.89	105.33	111.58
1	C	76	ILE	CB-CA-C	-5.84	102.12	110.83
1	B	110	ALA	CA-C-N	-5.80	114.78	120.52
1	B	110	ALA	C-N-CA	-5.80	114.78	120.52
1	D	297	ASP	N-CA-CB	-5.72	101.48	110.46
1	B	11	VAL	CB-CA-C	-5.67	104.44	111.08
1	B	346	ILE	CB-CA-C	5.67	116.96	111.23
1	A	148	SER	CA-C-N	-5.65	110.76	121.54
1	A	148	SER	C-N-CA	-5.65	110.76	121.54
1	B	347	LYS	N-CA-C	5.61	117.40	111.28
1	A	276	SER	N-CA-C	-5.58	105.30	111.71
1	C	110	ALA	CA-C-N	5.55	125.50	119.78
1	C	110	ALA	C-N-CA	5.55	125.50	119.78
1	C	19	THR	N-CA-C	-5.48	105.86	112.54
1	D	247	THR	N-CA-C	5.47	116.45	108.14
1	C	184	ASP	N-CA-C	-5.36	105.44	111.28
1	C	275	MET	N-CA-C	-5.35	105.36	111.14
1	C	146	VAL	CA-C-N	-5.28	114.82	120.47
1	C	146	VAL	C-N-CA	-5.28	114.82	120.47
1	D	379	VAL	N-CA-CB	5.25	115.40	109.99
1	B	16	ASP	N-CA-C	5.24	117.94	109.40
1	C	256	ILE	N-CA-C	5.24	115.97	110.62
1	C	297	ASP	N-CA-CB	-5.17	101.76	110.49
1	A	381	GLU	N-CA-C	-5.14	98.84	107.32
1	D	106	THR	CA-C-N	-5.13	116.83	123.19
1	D	106	THR	C-N-CA	-5.13	116.83	123.19
1	B	13	MET	CG-SD-CE	-5.11	89.66	100.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	364	ILE	CB-CA-C	-5.10	105.34	112.02
1	D	275	MET	N-CA-C	-5.07	105.88	111.71
1	D	367	GLY	CA-C-N	-5.04	116.94	123.19
1	D	367	GLY	C-N-CA	-5.04	116.94	123.19
1	A	100	ARG	CB-CA-C	-5.02	102.94	110.92
1	C	296	PRO	CA-C-N	5.00	131.10	121.54
1	C	296	PRO	C-N-CA	5.00	131.10	121.54

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	379	VAL	Peptide
1	B	380	PRO	Peptide
1	C	273	ASP	Peptide

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2917	0	2820	152	0
1	B	2859	0	2723	165	0
1	C	2884	0	2788	83	0
1	D	2930	0	2855	147	0
2	A	21	0	0	2	0
2	B	15	0	0	0	0
2	C	22	0	0	3	0
2	D	30	0	0	5	0
All	All	11678	0	11186	533	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 23.

All (533) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:379:VAL:HG12	1:A:380:PRO:CB	1.38	1.52
1:A:379:VAL:HG12	1:A:380:PRO:CG	1.10	1.52
1:A:379:VAL:CG1	1:A:380:PRO:CG	1.97	1.43
1:A:379:VAL:CG1	1:A:380:PRO:CB	2.00	1.38
1:A:379:VAL:CG1	1:A:380:PRO:HB3	1.54	1.36
1:B:13:MET:CE	1:B:60:ALA:HB1	1.58	1.32
1:A:379:VAL:CG1	1:A:380:PRO:HG3	1.58	1.29
1:A:355:LYS:HD3	1:A:398:PHE:CE1	1.68	1.28
1:D:273:ASP:O	1:D:274:VAL:HG23	1.16	1.25
1:B:241:ASN:C	1:B:242:LEU:HD23	1.66	1.21
1:A:118:LYS:HD3	1:A:378:GLU:OE2	1.36	1.21
1:B:13:MET:HE3	1:B:60:ALA:CB	1.74	1.17
1:B:115:ASN:HD22	1:B:115:ASN:N	1.37	1.15
1:A:355:LYS:HD3	1:A:398:PHE:CZ	1.82	1.13
1:A:379:VAL:HG13	1:A:380:PRO:HB3	1.16	1.12
1:B:13:MET:CE	1:B:60:ALA:CB	2.27	1.12
1:A:148:SER:O	1:A:149:ALA:CB	1.96	1.11
1:D:258:ARG:HH11	1:D:258:ARG:HG2	1.07	1.09
1:D:273:ASP:O	1:D:274:VAL:CG2	2.01	1.09
1:B:178:MET:HE3	1:B:180:MET:HE1	1.35	1.09
1:D:130:ILE:HG21	1:D:276:SER:HB2	1.29	1.09
1:A:148:SER:O	1:A:149:ALA:HB3	1.46	1.08
1:A:379:VAL:HG12	1:A:380:PRO:CD	1.84	1.07
1:B:379:VAL:O	1:B:379:VAL:HG13	1.48	1.07
1:C:130:ILE:HG21	1:C:276:SER:HB3	1.34	1.06
1:D:258:ARG:HH11	1:D:258:ARG:CG	1.69	1.05
1:B:115:ASN:H	1:B:115:ASN:ND2	1.34	1.02
1:A:379:VAL:HG12	1:A:380:PRO:HG3	1.17	1.02
1:B:242:LEU:HD23	1:B:242:LEU:N	1.62	1.01
1:A:202:ASN:OD1	1:A:240:LYS:NZ	1.95	1.00
1:D:363:THR:HG21	1:D:394:ILE:HA	1.44	0.99
1:D:107:VAL:HG22	1:D:293:LEU:CD1	1.92	0.97
1:A:274:VAL:O	1:A:277:ASP:HB2	1.66	0.96
1:D:107:VAL:HG22	1:D:293:LEU:HD13	1.48	0.95
1:A:355:LYS:CD	1:A:398:PHE:CE1	2.50	0.94
1:B:203:GLN:HG2	1:B:262:GLY:O	1.69	0.93
1:B:79:ASN:ND2	1:B:82:ARG:HG2	1.83	0.93
1:B:386:ASN:C	1:B:386:ASN:HD22	1.75	0.93
1:C:247:THR:HG21	1:C:252:ALA:HB2	1.50	0.93
1:B:380:PRO:HD2	1:B:380:PRO:O	1.69	0.93
1:B:379:VAL:O	1:B:379:VAL:CG1	2.16	0.91
1:C:258:ARG:HG2	1:C:258:ARG:HH11	1.33	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:258:ARG:HG2	1:D:258:ARG:NH1	1.85	0.89
1:A:363:THR:HG22	1:A:368:VAL:HG22	1.55	0.89
1:A:379:VAL:HG11	1:A:380:PRO:HG3	1.53	0.88
1:D:132:ARG:NH1	1:D:267:CYS:HB3	1.87	0.88
1:D:381:GLU:HB3	2:D:2028:HOH:O	1.73	0.88
1:A:188:ARG:HG2	1:A:228:ILE:HD11	1.55	0.88
1:C:208:SER:HB3	1:C:247:THR:HG23	1.54	0.88
1:B:105:GLY:HA3	1:B:293:LEU:HD11	1.57	0.87
1:B:158:SER:OG	1:B:160:ASN:ND2	2.07	0.86
1:B:149:ALA:O	1:B:171:PHE:O	1.93	0.86
1:B:79:ASN:H	1:B:79:ASN:HD22	1.20	0.86
1:D:160:ASN:ND2	1:D:160:ASN:H	1.70	0.86
1:A:380:PRO:HG2	1:A:382:LYS:HG3	1.58	0.84
1:B:363:THR:CG2	1:B:368:VAL:CG2	2.55	0.84
1:B:202:ASN:CG	1:B:240:LYS:HZ1	1.86	0.83
1:D:51:LYS:HB3	1:D:51:LYS:NZ	1.93	0.83
1:C:208:SER:HB3	1:C:247:THR:CG2	2.07	0.83
1:B:242:LEU:N	1:B:242:LEU:CD2	2.41	0.83
1:D:378:GLU:OE1	1:D:378:GLU:HA	1.76	0.83
1:C:386:ASN:HD22	1:C:386:ASN:C	1.87	0.82
1:B:13:MET:HE3	1:B:60:ALA:HB1	0.83	0.82
1:A:327:MET:HA	1:A:327:MET:HE2	1.61	0.82
1:C:267:CYS:HB2	2:C:2012:HOH:O	1.79	0.81
1:C:289:MET:HB3	1:C:305:ALA:HB3	1.62	0.81
1:A:327:MET:HA	1:A:327:MET:CE	2.10	0.81
1:B:178:MET:HE3	1:B:180:MET:CE	2.09	0.80
1:D:79:ASN:O	1:D:80:ALA:C	2.24	0.80
1:A:362:GLN:NE2	1:A:366:ASN:OD1	2.15	0.80
1:B:215:LYS:O	1:B:219:HIS:ND1	2.14	0.79
1:A:253:VAL:HG12	1:A:265:TRP:HH2	1.47	0.79
1:C:284:GLY:O	1:C:285:SER:HB2	1.82	0.79
1:D:379:VAL:O	1:D:382:LYS:HE2	1.84	0.78
1:A:355:LYS:CD	1:A:398:PHE:CZ	2.65	0.77
1:B:386:ASN:ND2	1:B:389:ASP:H	1.83	0.77
1:B:327:MET:HE3	1:B:360:SER:HB3	1.66	0.77
1:A:274:VAL:HG12	1:A:275:MET:N	1.98	0.77
1:C:273:ASP:HB2	1:C:275:MET:H	1.50	0.76
1:B:130:ILE:HD13	1:B:265:TRP:HB3	1.67	0.76
1:B:233:TYR:O	1:B:234:LYS:C	2.27	0.76
1:B:208:SER:OG	1:B:265:TRP:CH2	2.38	0.76
1:D:386:ASN:ND2	1:D:389:ASP:H	1.83	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:LYS:O	1:A:219:HIS:HB3	1.85	0.75
1:C:51:LYS:HB3	1:C:51:LYS:HZ2	1.51	0.75
1:C:355:LYS:HG3	1:C:401:MET:HE1	1.66	0.75
1:D:83:VAL:O	1:D:87:ASN:N	2.19	0.75
1:B:79:ASN:HD21	1:B:82:ARG:HG2	1.48	0.74
1:D:107:VAL:CG2	1:D:293:LEU:CD1	2.64	0.74
1:B:380:PRO:O	1:B:380:PRO:CD	2.35	0.74
1:D:362:GLN:HE21	1:D:366:ASN:HD21	1.34	0.74
1:B:363:THR:HG23	1:B:368:VAL:CG2	2.17	0.74
1:B:217:TYR:CE2	1:D:178:MET:HB2	2.23	0.73
1:B:274:VAL:O	1:B:277:ASP:N	2.20	0.73
1:D:21:ILE:HD12	1:D:21:ILE:H	1.53	0.73
1:D:114:VAL:HG22	1:D:117:ILE:HG13	1.71	0.73
1:A:3:LYS:HE2	1:A:32:GLU:O	1.88	0.73
1:D:386:ASN:C	1:D:386:ASN:HD22	1.96	0.73
1:B:241:ASN:C	1:B:242:LEU:CD2	2.56	0.73
1:B:384:ILE:HG13	1:B:384:ILE:O	1.86	0.73
1:B:348:GLU:OE1	1:B:348:GLU:N	2.21	0.73
1:D:132:ARG:CZ	1:D:267:CYS:HB3	2.19	0.72
1:D:104:ASP:O	1:D:104:ASP:OD1	2.07	0.72
1:D:291:SER:HB3	1:D:303:GLU:HG3	1.70	0.72
1:A:3:LYS:CE	1:A:32:GLU:O	2.38	0.72
1:C:284:GLY:O	1:C:285:SER:CB	2.39	0.71
1:D:51:LYS:HB3	1:D:51:LYS:HZ1	1.55	0.71
1:B:117:ILE:HG22	1:B:117:ILE:O	1.90	0.71
1:A:253:VAL:HG21	1:C:277:ASP:HB3	1.71	0.70
1:B:325:ASN:HB2	1:B:372:ASP:OD2	1.91	0.70
1:A:327:MET:HE1	1:A:360:SER:HB3	1.74	0.69
1:A:13:MET:HE1	1:A:64:ILE:HD11	1.74	0.69
1:C:386:ASN:ND2	1:C:389:ASP:H	1.90	0.69
1:C:247:THR:CG2	1:C:252:ALA:HB2	2.21	0.69
1:A:386:ASN:C	1:A:386:ASN:HD22	2.00	0.69
1:B:363:THR:HG22	1:B:368:VAL:CG2	2.22	0.69
1:D:146:VAL:CG2	1:D:175:GLY:H	2.07	0.69
1:B:208:SER:OG	1:B:265:TRP:HH2	1.74	0.68
1:D:51:LYS:NZ	1:D:51:LYS:CB	2.54	0.68
1:B:121:VAL:O	1:B:122:LYS:C	2.36	0.68
1:B:253:VAL:O	1:B:256:ILE:HG22	1.94	0.68
1:D:160:ASN:H	1:D:160:ASN:HD22	1.38	0.68
1:D:13:MET:HE3	1:D:72:LYS:HD2	1.76	0.68
1:A:188:ARG:HG2	1:A:228:ILE:CD1	2.23	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:379:VAL:H	1:A:380:PRO:HD3	1.58	0.68
1:B:178:MET:CE	1:B:180:MET:HE1	2.21	0.68
1:C:363:THR:HG21	1:C:394:ILE:HA	1.74	0.67
1:D:386:ASN:HD21	1:D:389:ASP:H	1.41	0.67
1:C:258:ARG:HH11	1:C:258:ARG:CG	2.08	0.67
1:A:251:ASP:O	1:A:255:ARG:HG2	1.95	0.67
1:A:379:VAL:HG12	1:A:380:PRO:N	2.04	0.66
1:A:160:ASN:HD21	1:A:162:GLU:HB2	1.60	0.66
1:D:200:ASP:OD2	1:D:339:LYS:NZ	2.17	0.66
1:A:362:GLN:HE21	1:A:366:ASN:CG	2.03	0.66
1:C:35:ILE:HD11	1:C:352:PHE:CE2	2.31	0.65
1:C:114:VAL:CG2	1:C:117:ILE:HG13	2.27	0.65
1:B:348:GLU:CD	1:B:348:GLU:H	2.03	0.65
1:D:81:GLN:O	1:D:84:GLU:HG2	1.96	0.65
1:D:362:GLN:HE21	1:D:366:ASN:ND2	1.93	0.65
1:A:274:VAL:O	1:A:277:ASP:N	2.30	0.65
1:D:359:ALA:O	1:D:363:THR:HG23	1.97	0.65
1:B:79:ASN:HD22	1:B:79:ASN:N	1.94	0.64
1:B:268:LYS:C	1:B:269:ASN:O	2.36	0.64
1:B:362:GLN:NE2	1:B:366:ASN:OD1	2.30	0.64
1:B:378:GLU:O	1:B:380:PRO:HG3	1.96	0.64
1:A:147:PRO:O	1:A:148:SER:HB3	1.97	0.64
1:D:82:ARG:O	1:D:83:VAL:C	2.39	0.64
1:D:154:LEU:C	1:D:154:LEU:HD13	2.22	0.64
1:A:115:ASN:N	1:A:115:ASN:HD22	1.96	0.64
1:B:240:LYS:O	1:B:242:LEU:HD21	1.97	0.64
1:D:258:ARG:HH11	1:D:258:ARG:CB	2.10	0.64
1:B:327:MET:CE	1:B:331:PHE:CE1	2.81	0.64
1:D:234:LYS:O	1:D:238:GLU:HG2	1.98	0.63
1:A:274:VAL:O	1:A:277:ASP:CB	2.43	0.63
1:D:377:SER:O	1:D:382:LYS:NZ	2.32	0.63
1:A:114:VAL:O	1:A:114:VAL:CG2	2.46	0.63
1:B:159:GLU:N	1:B:159:GLU:OE1	2.32	0.63
1:B:202:ASN:OD1	1:B:240:LYS:NZ	2.30	0.63
1:C:86:TYR:O	1:C:87:ASN:C	2.42	0.63
1:D:46:LEU:HD12	1:D:49:ARG:NH2	2.14	0.63
1:D:203:GLN:NE2	2:D:2014:HOH:O	2.19	0.63
1:A:295:SER:HB2	1:A:296:PRO:HD2	1.80	0.63
1:B:386:ASN:C	1:B:386:ASN:ND2	2.49	0.62
1:B:13:MET:HE1	1:B:60:ALA:CB	2.23	0.62
1:D:79:ASN:OD1	1:D:79:ASN:C	2.41	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:363:THR:CG2	1:A:368:VAL:HG22	2.28	0.62
1:D:114:VAL:CG2	1:D:117:ILE:HG13	2.29	0.62
1:B:368:VAL:HG23	1:B:368:VAL:O	1.98	0.62
1:C:218:ASP:OD1	1:C:268:LYS:HD3	2.00	0.62
1:C:355:LYS:CG	1:C:401:MET:HE1	2.28	0.62
1:C:226:GLN:O	1:C:230:GLU:HG2	2.00	0.62
1:D:130:ILE:HG21	1:D:276:SER:CB	2.19	0.62
1:D:180:MET:HE1	1:D:268:LYS:HE3	1.81	0.62
1:A:295:SER:HB2	1:A:296:PRO:CD	2.30	0.62
1:B:160:ASN:OD1	1:B:162:GLU:HG3	1.99	0.62
1:B:240:LYS:O	1:B:242:LEU:CD2	2.47	0.62
1:A:379:VAL:N	1:A:380:PRO:HD3	2.13	0.61
1:B:241:ASN:O	1:B:242:LEU:HD23	2.01	0.61
1:D:151:LYS:HE3	1:D:170:GLU:OE2	2.00	0.61
1:C:355:LYS:HG3	1:C:401:MET:CE	2.30	0.61
1:C:51:LYS:HB3	1:C:51:LYS:NZ	2.14	0.61
1:D:146:VAL:HG22	1:D:175:GLY:H	1.66	0.61
1:B:117:ILE:HD11	1:B:373:LEU:CD2	2.30	0.61
1:B:363:THR:HG23	1:B:368:VAL:HG21	1.82	0.61
1:D:258:ARG:CG	1:D:258:ARG:NH1	2.42	0.61
1:D:79:ASN:OD1	1:D:81:GLN:N	2.34	0.61
1:D:300:TYR:CE1	1:D:339:LYS:HE2	2.36	0.60
1:A:13:MET:CG	1:A:72:LYS:HB2	2.32	0.60
1:D:107:VAL:HG22	1:D:293:LEU:HD12	1.82	0.60
1:B:117:ILE:HD11	1:B:373:LEU:HD22	1.84	0.60
1:B:13:MET:CE	1:B:60:ALA:HB3	2.29	0.60
1:B:256:ILE:CG2	1:B:257:ILE:N	2.64	0.59
1:D:146:VAL:CG2	1:D:175:GLY:N	2.65	0.59
1:A:133:HIS:HB2	1:A:190:PHE:CD1	2.38	0.59
1:D:108:PHE:HB2	1:D:292:VAL:HG13	1.83	0.59
1:D:273:ASP:O	1:D:274:VAL:CB	2.49	0.59
1:A:160:ASN:C	1:A:160:ASN:HD22	2.10	0.59
1:B:249:ILE:O	1:B:253:VAL:HG13	2.03	0.59
1:C:63:ALA:HA	1:C:66:LYS:HD2	1.85	0.58
1:A:324:THR:O	1:A:370:THR:HG21	2.02	0.58
1:C:188:ARG:HG2	1:C:228:ILE:HD11	1.84	0.58
1:B:368:VAL:CG2	1:B:368:VAL:O	2.51	0.58
1:C:166:GLN:HG3	2:C:2013:HOH:O	2.03	0.58
1:C:258:ARG:HG2	1:C:258:ARG:NH1	2.08	0.58
1:D:252:ALA:O	1:D:256:ILE:HG12	2.03	0.58
1:B:217:TYR:HE2	1:D:178:MET:HB2	1.69	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:362:GLN:HE21	1:B:366:ASN:CG	2.11	0.58
1:A:115:ASN:HD22	1:A:115:ASN:H	1.51	0.58
1:A:379:VAL:HG12	1:A:380:PRO:CA	2.25	0.58
1:D:158:SER:OG	1:D:160:ASN:ND2	2.36	0.58
1:B:105:GLY:HA3	1:B:293:LEU:CD1	2.33	0.57
1:C:100:ARG:C	1:C:100:ARG:HH11	2.11	0.57
1:A:246:TYR:CD1	1:A:246:TYR:C	2.82	0.57
1:B:79:ASN:OD1	1:B:81:GLN:OE1	2.22	0.57
1:B:111:PRO:HG3	1:B:283:PHE:CD2	2.38	0.57
1:B:363:THR:CG2	1:B:368:VAL:HG22	2.35	0.57
1:D:78:PRO:HG2	1:D:92:TRP:O	2.05	0.57
1:A:386:ASN:ND2	1:A:389:ASP:H	2.02	0.57
1:C:359:ALA:O	1:C:363:THR:CG2	2.53	0.57
1:A:215:LYS:O	1:A:219:HIS:CB	2.53	0.56
1:D:141:ASN:CG	1:D:178:MET:HE3	2.31	0.56
1:A:219:HIS:CE1	1:A:223:ASP:OD2	2.59	0.56
1:C:386:ASN:HD21	1:C:389:ASP:H	1.53	0.56
1:A:32:GLU:N	1:A:33:PRO:CD	2.68	0.56
1:B:159:GLU:OE1	1:B:159:GLU:CA	2.53	0.56
1:B:200:ASP:C	1:B:200:ASP:OD1	2.49	0.56
1:B:272:GLY:O	1:B:276:SER:OG	2.24	0.56
1:B:364:ILE:HD13	1:B:373:LEU:HD13	1.88	0.56
1:A:370:THR:OG1	1:A:372:ASP:OD1	2.23	0.55
1:A:379:VAL:CB	1:A:380:PRO:CD	2.84	0.55
1:B:32:GLU:N	1:B:33:PRO:CD	2.68	0.55
1:D:235:GLU:CD	1:D:235:GLU:H	2.14	0.55
1:B:79:ASN:OD1	1:B:81:GLN:HB2	2.06	0.55
1:A:3:LYS:HE3	1:A:32:GLU:O	2.05	0.55
1:B:276:SER:O	1:B:280:ALA:HB2	2.07	0.55
1:C:160:ASN:C	1:C:160:ASN:HD22	2.13	0.55
1:A:160:ASN:C	1:A:160:ASN:ND2	2.64	0.55
1:A:364:ILE:HG23	1:A:369:MET:HE2	1.88	0.55
1:C:289:MET:HB3	1:C:305:ALA:CB	2.36	0.55
1:A:81:GLN:HB2	2:A:2009:HOH:O	2.06	0.55
1:C:158:SER:OG	1:C:160:ASN:ND2	2.40	0.55
1:D:13:MET:HG3	1:D:72:LYS:HB2	1.88	0.55
1:A:34:TYR:OH	1:A:399:GLU:OE2	2.21	0.55
1:D:18:MET:HA	1:D:21:ILE:HD13	1.89	0.55
1:D:80:ALA:O	1:D:84:GLU:HG2	2.07	0.55
1:B:216:THR:HG21	1:D:145:TYR:HB2	1.87	0.54
1:C:24:ARG:HG3	1:C:24:ARG:HH11	1.70	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:276:SER:O	1:B:280:ALA:CB	2.55	0.54
1:C:258:ARG:CG	1:C:258:ARG:NH1	2.66	0.54
1:A:363:THR:HG22	1:A:368:VAL:CG2	2.34	0.54
1:D:72:LYS:HE3	1:D:96:ASN:OD1	2.08	0.54
1:D:107:VAL:CG2	1:D:293:LEU:HD13	2.30	0.54
1:C:362:GLN:HE21	1:C:366:ASN:ND2	2.06	0.54
1:A:327:MET:HE3	1:A:330:ILE:HD12	1.89	0.54
1:A:304:ALA:HB1	1:A:328:ALA:HB1	1.89	0.54
1:D:21:ILE:HD12	1:D:21:ILE:N	2.21	0.54
1:A:216:THR:CG2	1:C:145:TYR:HB2	2.38	0.54
1:C:72:LYS:HB3	1:C:303:GLU:HG3	1.88	0.54
1:A:118:LYS:CD	1:A:378:GLU:OE2	2.31	0.53
1:B:219:HIS:C	1:B:219:HIS:CD2	2.86	0.53
1:C:35:ILE:HD11	1:C:352:PHE:CD2	2.43	0.53
1:A:379:VAL:CG1	1:A:380:PRO:CD	2.64	0.53
1:B:226:GLN:CA	1:B:226:GLN:OE1	2.56	0.53
1:D:269:ASN:O	1:D:273:ASP:OD1	2.26	0.53
1:B:79:ASN:ND2	1:B:79:ASN:H	1.96	0.53
1:C:100:ARG:NH1	1:C:101:ALA:HA	2.24	0.53
1:B:13:MET:HE1	1:B:60:ALA:HB3	1.90	0.53
1:B:325:ASN:CB	1:B:372:ASP:OD2	2.57	0.53
1:B:250:ASP:HB3	1:D:277:ASP:OD2	2.09	0.53
1:B:284:GLY:O	1:B:285:SER:CB	2.56	0.53
1:A:256:ILE:CG2	1:A:257:ILE:N	2.72	0.53
1:C:359:ALA:O	1:C:363:THR:HG23	2.09	0.53
1:D:13:MET:HB2	1:D:44:LEU:HD22	1.91	0.53
1:D:81:GLN:O	1:D:84:GLU:N	2.42	0.53
1:D:199:LEU:HD22	1:D:240:LYS:HE2	1.91	0.53
1:A:237:PHE:CD2	1:A:242:LEU:HD13	2.44	0.52
1:B:327:MET:CE	1:B:331:PHE:HE1	2.21	0.52
1:D:130:ILE:HD11	1:D:279:VAL:HG21	1.89	0.52
1:A:65:GLN:HG2	2:A:2020:HOH:O	2.08	0.52
1:B:79:ASN:HD22	1:B:82:ARG:HG2	1.70	0.52
1:B:133:HIS:HB2	1:B:190:PHE:CD1	2.44	0.52
1:A:380:PRO:HG2	1:A:382:LYS:CG	2.37	0.52
1:B:346:ILE:HG12	1:B:346:ILE:O	2.09	0.52
1:B:390:PHE:O	1:B:394:ILE:HG13	2.09	0.52
1:B:355:LYS:HE3	1:B:398:PHE:CZ	2.45	0.52
1:D:146:VAL:HG22	1:D:175:GLY:N	2.24	0.52
1:B:327:MET:HE2	1:B:331:PHE:CE1	2.44	0.52
1:D:226:GLN:O	1:D:230:GLU:HG2	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:166:GLN:HG3	2:D:2016:HOH:O	2.10	0.52
1:B:390:PHE:CE2	1:B:394:ILE:HD11	2.45	0.52
1:C:53:GLU:HG2	1:C:92:TRP:HH2	1.74	0.52
1:B:217:TYR:HB2	1:D:143:GLU:CD	2.35	0.52
1:A:386:ASN:C	1:A:386:ASN:ND2	2.65	0.51
1:B:208:SER:OG	1:B:265:TRP:CZ3	2.63	0.51
1:D:114:VAL:HG22	1:D:117:ILE:CG1	2.37	0.51
1:B:141:ASN:ND2	1:B:180:MET:HB3	2.25	0.51
1:D:7:LYS:HG3	1:D:346:ILE:HD11	1.93	0.51
1:B:115:ASN:N	1:B:115:ASN:ND2	2.11	0.51
1:B:241:ASN:O	1:B:242:LEU:CD2	2.58	0.51
1:C:13:MET:HB3	1:C:44:LEU:HD22	1.92	0.51
1:A:114:VAL:O	1:A:114:VAL:HG23	2.09	0.51
1:A:385:VAL:HB	1:A:389:ASP:HB3	1.93	0.51
1:D:154:LEU:HD12	2:D:2016:HOH:O	2.11	0.51
1:A:117:ILE:O	1:A:117:ILE:HG22	2.11	0.51
1:B:226:GLN:OE1	1:B:226:GLN:C	2.54	0.51
1:D:359:ALA:O	1:D:363:THR:CG2	2.59	0.51
1:A:127:PRO:O	1:A:262:GLY:HA2	2.11	0.51
1:B:143:GLU:OE1	1:D:216:THR:HB	2.10	0.51
1:D:132:ARG:HH12	1:D:267:CYS:HB3	1.74	0.51
1:A:379:VAL:H	1:A:380:PRO:CD	2.23	0.50
1:A:325:ASN:C	1:A:325:ASN:OD1	2.52	0.50
1:D:81:GLN:O	1:D:82:ARG:C	2.54	0.50
1:B:362:GLN:HE21	1:B:366:ASN:ND2	2.07	0.50
1:A:322:THR:O	1:A:386:ASN:HB2	2.11	0.50
1:B:160:ASN:H	1:B:160:ASN:HD22	1.59	0.50
1:B:226:GLN:OE1	1:B:226:GLN:O	2.30	0.50
1:C:325:ASN:ND2	1:C:328:ALA:H	2.10	0.50
1:D:108:PHE:CE2	1:D:131:ALA:HB2	2.46	0.50
1:D:258:ARG:NH1	1:D:258:ARG:CB	2.73	0.50
1:A:211:ASP:HB2	1:A:218:ASP:HB3	1.94	0.50
1:A:385:VAL:HB	1:A:389:ASP:CB	2.42	0.50
1:A:143:GLU:OE1	1:C:216:THR:HB	2.12	0.50
1:A:253:VAL:O	1:A:256:ILE:HG22	2.12	0.50
1:D:108:PHE:CD2	1:D:131:ALA:HB2	2.46	0.50
1:D:295:SER:HB2	1:D:296:PRO:HD2	1.93	0.50
1:A:188:ARG:CG	1:A:228:ILE:HD11	2.33	0.50
1:B:130:ILE:HD13	1:B:265:TRP:CB	2.37	0.50
1:B:149:ALA:O	1:B:171:PHE:HB2	2.12	0.50
1:C:386:ASN:C	1:C:386:ASN:ND2	2.60	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:146:VAL:HG23	1:D:175:GLY:H	1.75	0.50
1:B:295:SER:HB2	1:B:296:PRO:CD	2.42	0.49
1:A:355:LYS:CE	1:A:398:PHE:CZ	2.96	0.49
1:B:355:LYS:HD2	1:B:401:MET:HE3	1.93	0.49
1:C:143:GLU:HG2	1:C:178:MET:HG2	1.94	0.49
1:A:386:ASN:HD21	1:A:389:ASP:H	1.58	0.49
1:C:100:ARG:NH1	1:C:100:ARG:O	2.45	0.49
1:D:46:LEU:CD1	1:D:49:ARG:NH2	2.75	0.49
1:A:216:THR:HG21	1:C:145:TYR:HB2	1.94	0.49
1:B:327:MET:CE	1:B:360:SER:HB3	2.40	0.49
1:B:400:GLY:O	1:B:401:MET:O	2.30	0.49
1:A:91:MET:O	1:A:92:TRP:C	2.55	0.49
1:B:274:VAL:O	1:B:277:ASP:HB2	2.11	0.49
1:C:82:ARG:O	1:C:83:VAL:C	2.55	0.49
1:A:256:ILE:HG23	1:A:257:ILE:N	2.27	0.49
1:C:114:VAL:HG22	1:C:117:ILE:HG13	1.95	0.49
1:A:180:MET:HG2	1:A:181:HIS:N	2.28	0.49
1:C:397:THR:O	1:C:401:MET:HG3	2.13	0.49
1:D:146:VAL:HG21	1:D:171:PHE:CD2	2.48	0.49
1:D:160:ASN:ND2	1:D:160:ASN:N	2.48	0.49
1:D:108:PHE:CE2	1:D:131:ALA:CB	2.95	0.48
1:A:195:PHE:O	1:A:196:ASN:C	2.56	0.48
1:B:363:THR:HG22	1:B:368:VAL:HG23	1.94	0.48
1:B:91:MET:HB2	1:B:91:MET:HE2	1.52	0.48
1:D:368:VAL:HG11	1:D:393:GLU:HG3	1.96	0.48
1:A:355:LYS:CE	1:A:398:PHE:CE1	2.96	0.48
1:D:89:LYS:O	1:D:90:LYS:HG3	2.13	0.48
1:A:211:ASP:OD1	1:A:211:ASP:C	2.56	0.48
1:B:352:PHE:O	1:B:353:ALA:C	2.56	0.48
1:D:378:GLU:OE1	1:D:378:GLU:CA	2.55	0.48
1:A:158:SER:OG	1:A:160:ASN:ND2	2.46	0.48
1:A:379:VAL:CB	1:A:380:PRO:HD3	2.43	0.48
1:A:398:PHE:O	1:A:400:GLY:N	2.44	0.48
1:D:249:ILE:O	1:D:253:VAL:HG13	2.14	0.48
1:D:66:LYS:NZ	2:D:2009:HOH:O	2.47	0.47
1:A:352:PHE:CD1	1:A:352:PHE:C	2.92	0.47
1:A:362:GLN:HE21	1:A:366:ASN:ND2	2.12	0.47
1:B:237:PHE:CD2	1:B:242:LEU:HB2	2.49	0.47
1:C:160:ASN:C	1:C:160:ASN:ND2	2.70	0.47
1:A:237:PHE:HA	1:A:242:LEU:HD11	1.96	0.47
1:B:78:PRO:HA	1:B:82:ARG:HG3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:149:ALA:O	1:B:171:PHE:C	2.56	0.47
1:B:68:GLY:C	1:B:69:VAL:HG23	2.39	0.47
1:B:240:LYS:HB3	1:B:240:LYS:HE2	1.43	0.47
1:D:246:TYR:CD2	1:D:246:TYR:C	2.93	0.47
1:B:246:TYR:CD1	1:B:247:THR:N	2.83	0.47
1:D:238:GLU:O	1:D:239:ALA:C	2.58	0.47
1:A:13:MET:HA	1:A:42:TYR:O	2.14	0.47
1:A:220:ARG:HA	1:A:220:ARG:CZ	2.45	0.47
1:B:355:LYS:CE	1:B:398:PHE:CZ	2.98	0.47
1:D:246:TYR:O	1:D:247:THR:HB	2.15	0.47
1:B:149:ALA:HB3	1:D:159:GLU:CG	2.45	0.47
1:A:209:THR:OG1	1:A:211:ASP:HB3	2.15	0.46
1:A:379:VAL:CB	1:A:380:PRO:HG3	2.37	0.46
1:B:83:VAL:HA	1:B:88:LEU:HG	1.97	0.46
1:B:112:ILE:HB	1:B:288:MET:HE3	1.96	0.46
1:A:327:MET:HE2	1:A:327:MET:CA	2.41	0.46
1:C:17:GLU:O	1:C:20:ARG:HB3	2.16	0.46
1:D:199:LEU:CD2	1:D:240:LYS:HE2	2.44	0.46
1:B:246:TYR:CD1	1:B:246:TYR:C	2.93	0.46
1:B:65:GLN:O	1:B:299:LYS:HE3	2.16	0.46
1:B:233:TYR:O	1:B:235:GLU:N	2.48	0.46
1:A:160:ASN:ND2	1:A:162:GLU:H	2.14	0.46
1:A:211:ASP:O	1:A:211:ASP:CG	2.58	0.46
1:B:114:VAL:HG22	1:B:117:ILE:HB	1.96	0.46
1:C:48:ASN:OD1	1:C:51:LYS:NZ	2.40	0.46
1:D:104:ASP:OD1	1:D:104:ASP:C	2.59	0.46
1:B:35:ILE:HD11	1:B:352:PHE:CE2	2.51	0.46
1:B:115:ASN:HD22	1:B:115:ASN:H	0.61	0.46
1:A:379:VAL:CB	1:A:380:PRO:CG	2.86	0.45
1:B:256:ILE:HG22	1:B:257:ILE:N	2.31	0.45
1:B:363:THR:HG22	1:B:368:VAL:HG22	1.94	0.45
1:D:37:LEU:HD12	1:D:37:LEU:HA	1.40	0.45
1:D:72:LYS:CE	1:D:96:ASN:OD1	2.63	0.45
1:A:363:THR:CG2	1:A:368:VAL:CG2	2.94	0.45
1:C:235:GLU:N	1:C:235:GLU:CD	2.75	0.45
1:C:362:GLN:HE21	1:C:366:ASN:HD21	1.62	0.45
1:A:126:LYS:O	1:A:127:PRO:C	2.57	0.45
1:C:258:ARG:HD3	2:C:2020:HOH:O	2.15	0.45
1:A:338:LYS:HE3	1:A:338:LYS:HB2	1.30	0.45
1:B:247:THR:OG1	1:B:248:LEU:N	2.50	0.45
1:A:371:LYS:O	1:A:374:ALA:HB3	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:68:GLY:C	1:B:69:VAL:CG2	2.90	0.45
1:B:356:LEU:O	1:B:357:GLU:C	2.58	0.45
1:A:323:SER:OG	1:A:371:LYS:CB	2.65	0.45
1:A:355:LYS:HE2	1:A:398:PHE:CE2	2.52	0.45
1:B:327:MET:HE3	1:B:360:SER:CB	2.43	0.45
1:C:208:SER:HA	1:C:247:THR:O	2.16	0.45
1:A:104:ASP:CG	1:A:104:ASP:O	2.58	0.45
1:A:293:LEU:HB3	1:A:301:GLU:HB3	1.99	0.45
1:C:224:ILE:O	1:C:228:ILE:HG12	2.16	0.45
1:A:155:VAL:HG21	1:C:155:VAL:HG21	1.98	0.45
1:B:216:THR:HA	1:B:219:HIS:CE1	2.52	0.45
1:B:275:MET:HA	1:B:278:MET:HB3	1.98	0.45
1:D:133:HIS:O	1:D:134:ALA:HB2	2.17	0.45
1:D:386:ASN:ND2	1:D:386:ASN:C	2.70	0.45
1:A:369:MET:HE2	1:A:369:MET:HB3	1.88	0.45
1:B:130:ILE:CD1	1:B:265:TRP:HB3	2.43	0.45
1:C:354:THR:HG22	1:C:358:GLN:NE2	2.32	0.45
1:A:23:TRP:HH2	1:A:71:VAL:HG22	1.83	0.44
1:B:355:LYS:HD2	1:B:401:MET:CE	2.47	0.44
1:D:197:TYR:CE1	1:D:292:VAL:HG11	2.52	0.44
1:C:77:THR:HA	1:C:78:PRO:HD3	1.76	0.44
1:C:295:SER:OG	1:C:297:ASP:HB3	2.18	0.44
1:D:203:GLN:HG3	1:D:264:VAL:HG23	2.00	0.44
1:D:268:LYS:HE2	1:D:268:LYS:HB3	1.60	0.44
1:D:132:ARG:HG3	1:D:135:TYR:CD1	2.52	0.44
1:D:26:ILE:O	1:D:27:LYS:C	2.59	0.44
1:D:379:VAL:HA	1:D:380:PRO:HD3	1.65	0.44
1:A:13:MET:HG3	1:A:72:LYS:HB2	1.99	0.44
1:A:368:VAL:HA	1:A:383:LYS:O	2.17	0.44
1:D:247:THR:HG23	1:D:252:ALA:HB2	1.99	0.44
1:A:116:SER:CB	1:A:379:VAL:HG23	2.48	0.44
1:A:99:ILE:HG21	1:A:301:GLU:OE2	2.18	0.44
1:D:180:MET:HE2	1:D:217:TYR:OH	2.18	0.44
1:A:7:LYS:HB2	1:A:346:ILE:HD11	1.99	0.43
1:A:100:ARG:O	1:A:104:ASP:N	2.50	0.43
1:B:240:LYS:O	1:B:242:LEU:HD23	2.16	0.43
1:B:363:THR:CG2	1:B:368:VAL:HG23	2.45	0.43
1:D:228:ILE:O	1:D:232:GLU:HB2	2.17	0.43
1:A:116:SER:HB3	1:A:379:VAL:HG23	1.99	0.43
1:C:95:PRO:O	1:C:99:ILE:HD12	2.17	0.43
1:A:209:THR:HB	1:A:218:ASP:CG	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:344:ASP:O	1:B:345:GLY:C	2.61	0.43
1:B:368:VAL:HG21	1:B:393:GLU:HG2	2.00	0.43
1:C:379:VAL:HA	1:C:380:PRO:HD3	1.53	0.43
1:B:363:THR:HG23	1:B:368:VAL:HG22	1.97	0.43
1:A:327:MET:HA	1:A:327:MET:HE3	1.97	0.43
1:A:380:PRO:HB2	1:A:381:GLU:H	1.11	0.43
1:B:201:MET:HE3	1:B:201:MET:HB3	1.71	0.43
1:B:202:ASN:CG	1:B:240:LYS:NZ	2.67	0.43
1:B:400:GLY:O	1:B:401:MET:C	2.61	0.43
1:C:355:LYS:CG	1:C:401:MET:CE	2.93	0.43
1:D:295:SER:OG	1:D:299:LYS:HB2	2.18	0.43
1:A:233:TYR:O	1:A:234:LYS:C	2.61	0.43
1:D:211:ASP:OD1	1:D:211:ASP:C	2.60	0.43
1:B:325:ASN:OD1	1:B:325:ASN:C	2.61	0.42
1:B:398:PHE:O	1:B:400:GLY:O	2.36	0.42
1:C:202:ASN:HD21	1:C:240:LYS:HD3	1.85	0.42
1:D:21:ILE:H	1:D:21:ILE:CD1	2.28	0.42
1:A:15:GLY:O	1:A:20:ARG:HD2	2.18	0.42
1:B:278:MET:O	1:B:282:ALA:HB2	2.19	0.42
1:D:154:LEU:C	1:D:154:LEU:CD1	2.91	0.42
1:C:202:ASN:ND2	1:C:240:LYS:HD3	2.34	0.42
1:D:13:MET:HG3	1:D:72:LYS:CB	2.49	0.42
1:D:295:SER:C	1:D:297:ASP:H	2.27	0.42
1:A:13:MET:HG2	1:A:72:LYS:HB2	2.01	0.42
1:D:52:THR:O	1:D:53:GLU:C	2.60	0.42
1:D:184:ASP:OD1	1:D:220:ARG:HD3	2.20	0.42
1:A:242:LEU:HD12	1:A:242:LEU:O	2.19	0.42
1:D:100:ARG:HG3	1:D:293:LEU:HD11	2.02	0.42
1:B:22:ILE:HG21	1:B:326:SER:HB3	2.02	0.42
1:B:179:GLY:HA2	1:D:178:MET:O	2.19	0.42
1:C:268:LYS:O	1:C:269:ASN:C	2.61	0.42
1:C:275:MET:O	1:C:278:MET:HB3	2.19	0.42
1:D:77:THR:HA	1:D:78:PRO:HD2	1.60	0.42
1:B:81:GLN:CD	1:B:81:GLN:H	2.27	0.42
1:A:7:LYS:HB3	1:A:8:VAL:H	1.58	0.42
1:A:208:SER:HA	1:A:247:THR:O	2.20	0.42
1:C:208:SER:HB3	1:C:247:THR:HG22	1.93	0.42
1:A:17:GLU:O	1:A:21:ILE:HD12	2.20	0.41
1:B:11:VAL:O	1:B:70:GLY:HA2	2.19	0.41
1:C:350:VAL:CG2	1:C:351:ASP:N	2.82	0.41
1:D:146:VAL:HG22	1:D:175:GLY:C	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:268:LYS:O	1:D:269:ASN:C	2.63	0.41
1:A:96:ASN:OD1	1:A:303:GLU:OE2	2.38	0.41
1:B:226:GLN:OE1	1:B:226:GLN:HA	2.20	0.41
1:B:361:VAL:CG1	1:B:365:GLU:OE2	2.69	0.41
1:A:115:ASN:N	1:A:115:ASN:ND2	2.67	0.41
1:D:295:SER:HB2	1:D:296:PRO:CD	2.51	0.41
1:A:18:MET:O	1:A:21:ILE:HB	2.20	0.41
1:A:114:VAL:HG22	1:A:117:ILE:HB	2.02	0.41
1:D:295:SER:C	1:D:297:ASP:N	2.78	0.41
1:C:24:ARG:HG3	1:C:24:ARG:NH1	2.34	0.41
1:C:114:VAL:HG22	1:C:117:ILE:CG1	2.51	0.41
1:D:32:GLU:N	1:D:33:PRO:CD	2.83	0.41
1:D:211:ASP:O	1:D:211:ASP:CG	2.61	0.41
1:D:280:ALA:HA	1:D:289:MET:HE1	2.02	0.41
1:A:108:PHE:CE2	1:A:131:ALA:HB2	2.56	0.41
1:B:111:PRO:HG3	1:B:283:PHE:CE2	2.56	0.41
1:B:355:LYS:HE3	1:B:398:PHE:CE2	2.54	0.41
1:C:268:LYS:HB3	1:C:268:LYS:HE2	1.79	0.41
1:D:133:HIS:HB2	1:D:190:PHE:CD1	2.56	0.41
1:D:154:LEU:HB2	1:D:168:ILE:HD11	2.03	0.41
1:A:160:ASN:HD22	1:A:161:GLY:N	2.18	0.41
1:A:348:GLU:H	1:A:348:GLU:CD	2.28	0.41
1:A:380:PRO:CG	1:A:381:GLU:N	2.74	0.41
1:B:216:THR:CG2	1:D:145:TYR:HB2	2.51	0.41
1:A:23:TRP:CE3	1:A:73:CYS:HB2	2.56	0.41
1:A:245:PHE:HE2	1:A:255:ARG:HD2	1.87	0.41
1:D:82:ARG:O	1:D:86:TYR:N	2.39	0.41
1:D:141:ASN:OD1	1:D:141:ASN:C	2.64	0.41
1:A:384:ILE:O	1:A:384:ILE:HG13	2.18	0.40
1:C:13:MET:HE1	1:C:301:GLU:HG3	2.03	0.40
1:D:230:GLU:HG2	1:D:230:GLU:H	1.63	0.40
1:D:289:MET:HB3	1:D:305:ALA:CB	2.51	0.40
1:B:13:MET:HA	1:B:42:TYR:O	2.21	0.40
1:B:208:SER:HB3	1:B:247:THR:O	2.22	0.40
1:C:160:ASN:HD22	1:C:161:GLY:N	2.19	0.40
1:C:359:ALA:O	1:C:363:THR:HG22	2.21	0.40
1:D:77:THR:OG1	1:D:94:SER:HB2	2.22	0.40
1:D:151:LYS:CE	1:D:170:GLU:OE2	2.69	0.40
1:B:394:ILE:O	1:B:395:ARG:C	2.64	0.40
1:C:300:TYR:CE1	1:C:339:LYS:HE2	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	368/402 (92%)	338 (92%)	24 (6%)	6 (2%)	7	6
1	B	364/402 (90%)	329 (90%)	28 (8%)	7 (2%)	6	5
1	C	363/402 (90%)	342 (94%)	16 (4%)	5 (1%)	9	7
1	D	367/402 (91%)	347 (95%)	17 (5%)	3 (1%)	16	17
All	All	1462/1608 (91%)	1356 (93%)	85 (6%)	21 (1%)	9	7

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	148	SER
1	A	380	PRO
1	B	122	LYS
1	B	150	GLY
1	B	285	SER
1	B	380	PRO
1	C	274	VAL
1	C	285	SER
1	D	274	VAL
1	A	149	ALA
1	C	297	ASP
1	D	83	VAL
1	A	379	VAL
1	B	399	GLU
1	A	399	GLU
1	B	269	ASN
1	B	378	GLU
1	C	87	ASN
1	D	136	GLY
1	A	400	GLY
1	C	83	VAL

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	302/347 (87%)	265 (88%)	37 (12%)	4	4
1	B	291/347 (84%)	247 (85%)	44 (15%)	3	2
1	C	299/347 (86%)	273 (91%)	26 (9%)	9	10
1	D	306/347 (88%)	267 (87%)	39 (13%)	4	4
All	All	1198/1388 (86%)	1052 (88%)	146 (12%)	5	4

All (146) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LYS
1	A	8	VAL
1	A	18	MET
1	A	37	LEU
1	A	53	GLU
1	A	100	ARG
1	A	113	VAL
1	A	114	VAL
1	A	115	ASN
1	A	118	LYS
1	A	125	LYS
1	A	159	GLU
1	A	160	ASN
1	A	162	GLU
1	A	164	SER
1	A	205	LEU
1	A	210	LYS
1	A	211	ASP
1	A	220	ARG
1	A	242	LEU
1	A	255	ARG
1	A	256	ILE
1	A	269	ASN
1	A	274	VAL

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Mol	Chain	Res	Type
1	A	276	SER
1	A	292	VAL
1	A	322	THR
1	A	327	MET
1	A	329	THR
1	A	338	LYS
1	A	368	VAL
1	A	377	SER
1	A	379	VAL
1	A	384	ILE
1	A	386	ASN
1	A	393	GLU
1	A	401	MET
1	B	21	ILE
1	B	25	LEU
1	B	37	LEU
1	B	47	GLU
1	B	53	GLU
1	B	79	ASN
1	B	83	VAL
1	B	87	ASN
1	B	91	MET
1	B	94	SER
1	B	114	VAL
1	B	115	ASN
1	B	117	ILE
1	B	159	GLU
1	B	160	ASN
1	B	164	SER
1	B	172	ASP
1	B	186	SER
1	B	200	ASP
1	B	201	MET
1	B	205	LEU
1	B	216	THR
1	B	219	HIS
1	B	226	GLN
1	B	230	GLU
1	B	240	LYS
1	B	242	LEU
1	B	256	ILE
1	B	257	ILE

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Mol	Chain	Res	Type
1	B	265	TRP
1	B	271	ASP
1	B	276	SER
1	B	292	VAL
1	B	322	THR
1	B	326	SER
1	B	329	THR
1	B	346	ILE
1	B	348	GLU
1	B	379	VAL
1	B	380	PRO
1	B	384	ILE
1	B	386	ASN
1	B	393	GLU
1	B	401	MET
1	C	21	ILE
1	C	51	LYS
1	C	91	MET
1	C	100	ARG
1	C	114	VAL
1	C	146	VAL
1	C	160	ASN
1	C	165	ARG
1	C	199	LEU
1	C	200	ASP
1	C	205	LEU
1	C	211	ASP
1	C	212	THR
1	C	230	GLU
1	C	253	VAL
1	C	257	ILE
1	C	258	ARG
1	C	267	CYS
1	C	268	LYS
1	C	273	ASP
1	C	289	MET
1	C	297	ASP
1	C	350	VAL
1	C	363	THR
1	C	379	VAL
1	C	386	ASN
1	D	2	SER

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Mol	Chain	Res	Type
1	D	13	MET
1	D	37	LEU
1	D	46	LEU
1	D	62	ARG
1	D	76	ILE
1	D	77	THR
1	D	79	ASN
1	D	84	GLU
1	D	93	LYS
1	D	94	SER
1	D	114	VAL
1	D	142	VAL
1	D	160	ASN
1	D	164	SER
1	D	202	ASN
1	D	205	LEU
1	D	211	ASP
1	D	220	ARG
1	D	230	GLU
1	D	235	GLU
1	D	238	GLU
1	D	253	VAL
1	D	257	ILE
1	D	258	ARG
1	D	268	LYS
1	D	273	ASP
1	D	276	SER
1	D	285	SER
1	D	288	MET
1	D	289	MET
1	D	299	LYS
1	D	301	GLU
1	D	325	ASN
1	D	350	VAL
1	D	363	THR
1	D	378	GLU
1	D	379	VAL
1	D	386	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	115	ASN
1	A	160	ASN
1	A	219	HIS
1	A	241	ASN
1	A	362	GLN
1	A	386	ASN
1	B	65	GLN
1	B	79	ASN
1	B	87	ASN
1	B	96	ASN
1	B	115	ASN
1	B	141	ASN
1	B	160	ASN
1	B	203	GLN
1	B	362	GLN
1	B	386	ASN
1	C	96	ASN
1	C	160	ASN
1	C	181	HIS
1	C	226	GLN
1	C	325	ASN
1	C	358	GLN
1	C	366	ASN
1	C	386	ASN
1	D	48	ASN
1	D	160	ASN
1	D	231	ASN
1	D	325	ASN
1	D	366	ASN
1	D	386	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	376/402 (93%)	0.21	17 (4%) 38 44	31, 50, 80, 121	0
1	B	372/402 (92%)	0.29	16 (4%) 40 45	33, 53, 86, 104	0
1	C	371/402 (92%)	-0.09	4 (1%) 78 82	26, 42, 75, 97	0
1	D	375/402 (93%)	-0.03	14 (3%) 45 51	27, 44, 75, 100	0
All	All	1494/1608 (92%)	0.10	51 (3%) 48 55	26, 47, 81, 121	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	211	ASP	4.7
1	A	213	ILE	4.4
1	B	216	THR	4.4
1	B	217	TYR	4.1
1	B	270	TYR	3.8
1	D	135	TYR	3.7
1	A	212	THR	3.7
1	A	322	THR	3.7
1	C	211	ASP	3.4
1	A	401	MET	3.3
1	A	217	TYR	3.3
1	D	211	ASP	3.1
1	B	379	VAL	3.1
1	B	245	PHE	3.1
1	A	271	ASP	3.0
1	D	274	VAL	3.0
1	B	215	LYS	2.9
1	C	274	VAL	2.8
1	D	136	GLY	2.6
1	B	265	TRP	2.6
1	A	272	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	216	THR	2.6
1	D	78	PRO	2.6
1	C	80	ALA	2.6
1	D	137	ASP	2.6
1	B	209	THR	2.6
1	D	86	TYR	2.6
1	A	400	GLY	2.5
1	A	379	VAL	2.5
1	B	271	ASP	2.4
1	D	269	ASN	2.4
1	B	269	ASN	2.3
1	A	387	THR	2.3
1	A	269	ASN	2.3
1	D	273	ASP	2.3
1	D	91	MET	2.3
1	C	78	PRO	2.2
1	B	149	ALA	2.2
1	B	381	GLU	2.2
1	D	134	ALA	2.2
1	A	214	SER	2.2
1	B	378	GLU	2.2
1	D	80	ALA	2.2
1	A	376	LEU	2.2
1	A	18	MET	2.2
1	D	172	ASP	2.1
1	D	85	GLU	2.1
1	A	381	GLU	2.1
1	B	274	VAL	2.1
1	B	234	LYS	2.0
1	B	380	PRO	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.